

Exercise 03 Raphael Michel and Florian Stoertz

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1 Linear Programs

1.1 Implementation

Using the definition given in the scipy manal, we write our graphical model as a linear program:

$$\min \quad c^T \mu \quad (1)$$

$$\text{s.t.} \quad A_u \mu = b_u \quad (2)$$

$$A_e \mu = b_e \quad (3)$$

For the given graphical model, we have

$$\mu^T = [\mu_0(0, z_0), \mu_0(1, z_0), \mu_1(0, z_1), \mu_1(1, z_1), \mu_2(0, z_2), \mu_2(1, z_2), \quad (4)$$

$$\mu_{01}(0, 0, z_0, z_1), \mu_{01}(0, 1, z_0, z_1), \mu_{01}(1, 0, z_0, z_1), \mu_{01}(1, 1, z_0, z_1), \quad (5)$$

$$\mu_{02}(0, 0, z_0, z_2), \mu_{02}(0, 1, z_0, z_2), \mu_{02}(1, 0, z_0, z_2), \mu_{02}(1, 1, z_0, z_2), \quad (6)$$

$$\mu_{12}(0, 0, z_1, z_2), \mu_{12}(0, 1, z_1, z_2), \mu_{12}(1, 0, z_1, z_2), \mu_{12}(1, 1, z_1, z_2)] \quad (7)$$

and a cost vector of

$$c^T = [\psi_0(0), \psi_0(1), \psi_1(0), \psi_1(1), \psi_2(0), \psi_2(1), \quad (8)$$

$$\psi_p(0, 0), \psi_p(0, 1), \psi_p(1, 0), \psi_p(1, 1), \quad (9)$$

$$\psi_p(0, 0), \psi_p(0, 1), \psi_p(1, 0), \psi_p(1, 1), \quad (10)$$

$$\psi_p(0, 0), \psi_p(0, 1), \psi_p(1, 0), \psi_p(1, 1)]. \quad (11)$$

From our constraint $\mu_{ij}(k, l) \geq 0$ we get the inequality constraint

$$\begin{bmatrix} 1 & & & \\ & 1 & & \\ & & \ddots & \\ & & & 1 \end{bmatrix} \mu \geq \begin{bmatrix} 0 \\ \vdots \\ 0 \end{bmatrix}. \quad (12)$$

For the equality constraints, we get

$$\left[\begin{array}{c|ccc} C & & & 0 \\ \hline & E & & \\ D & & E & \\ & & & E \end{array} \right] \mu = \left[\begin{array}{c} \mathbb{1}_{3 \times 1} \\ 0 \end{array} \right]. \quad (13)$$

In this block matrix, the C part realizes the condition $\sum_k \mu_i(k) = 1$:

$$C = \begin{bmatrix} 1 & 1 & & & \\ & & 1 & 1 & \\ & & & & 1 & 1 \end{bmatrix} \quad (14)$$

And the lower part realizes the other two conditions $\sum_k \mu_{ij}(k, l) = \mu_j(l)$ and $\sum_l \mu_{ij}(k, l) = \mu_i(k)$:

$$D = \begin{bmatrix} -1 & & & & & & \\ & -1 & & & & & \\ & & -1 & & & & \\ & & & -1 & & & \\ -1 & & & & & & \\ & -1 & & & & & \\ & & -1 & & & & \\ & & & -1 & & & \\ & & & & -1 & & \\ & & & & & -1 & \\ & & & & & & -1 \end{bmatrix} \quad (15)$$

$$E = \begin{bmatrix} 1 & 1 & & & \\ & & 1 & 1 & \\ 1 & & 1 & & \\ & 1 & & 1 & \end{bmatrix} \quad (16)$$

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In [28]: import numpy as np
         from scipy.linalg import block_diag
         from scipy.optimize import linprog
```

```
In [35]: def solve(beta):
         unary00 = .1
         unary01 = .1
         unary10 = .1
         unary11 = .9
         unary20 = .9
         unary21 = .1
         c = np.array([
             unary00, unary01, unary10, unary11, unary20, unary21,
             0, beta, beta, 0,
             0, beta, beta, 0,
             0, beta, beta, 0,
```

```

])
A_ub = - np.eye(18)
b_ub = np.zeros(18)

C = np.array([
    [1, 1, 0, 0, 0, 0],
    [0, 0, 1, 1, 0, 0],
    [0, 0, 0, 0, 1, 1]
])
D = np.array([
    [-1, 0, 0, 0, 0, 0],
    [0, -1, 0, 0, 0, 0],
    [0, 0, -1, 0, 0, 0],
    [0, 0, 0, -1, 0, 0],
    [-1, 0, 0, 0, 0, 0],
    [0, -1, 0, 0, 0, 0],
    [0, 0, 0, 0, -1, 0],
    [0, 0, 0, 0, 0, -1],
    [0, 0, -1, 0, 0, 0],
    [0, 0, 0, -1, 0, 0],
    [0, 0, 0, 0, -1, 0],
    [0, 0, 0, 0, 0, -1],
])
E = np.array([
    [1, 1, 0, 0],
    [0, 0, 1, 1],
    [1, 0, 1, 0],
    [0, 1, 0, 1]
])
F = block_diag(E, E, E)
A_eq = np.vstack((
    np.hstack((C, np.zeros((3, 12)))),
    np.hstack((D, F))
))
b_eq = np.hstack((np.ones(3), np.zeros(12)))
x = linprog(c, A_ub, b_ub, A_eq, b_eq)
return x

```

In [38]: *# Solving for an attractive potential*
solve(1.0)

```

Out[38]:      fun: 1.1000000000000001
      message: 'Optimization terminated successfully.'
      nit: 17
      slack: array([ 1.,  0.,  1.,  0.,  1.,  0.,  1.,  0.,  0.,  0.,  1.,  0.,  0.,
                    0.,  1.,  0.,  0.,  0.])
      status: 0
      success: True

```

```
x: array([ 1.,  0.,  1.,  0.,  1.,  0.,  1.,  0.,  0.,  0.,  1.,  0.,  0.,
          0.,  1.,  0.,  0.,  0.] )
```

For the attractive potential we get a state with the minimal energy of 1.1.

```
In [37]: solve(-1.0)
```

```
Out[37]:      fun: -1.8999999999999999
      message: 'Optimization terminated successfully.'
      nit: 17
      slack: array([ 0.5,  0.5,  0.5,  0.5,  0.5,  0.5,  0. ,  0.5,  0.5,  0. ,  0. ,
                    0.5,  0.5,  0. ,  0. ,  0.5,  0.5,  0. ])
      status: 0
      success: True
      x: array([ 0.5,  0.5,  0.5,  0.5,  0.5,  0.5,  0. ,  0.5,  0.5,  0. ,  0. ,
                0.5,  0.5,  0. ,  0. ,  0.5,  0.5,  0. ] )
```

For the repulsive potential we get a state with a negative minimal energy. This is not a surprise, as we posed the attractiveness as a requirement for this method in the lecture.